LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
Data File : BP021816.D
Acq On : 04 Sep 2024 12:38
Operator : RC/JU
Sample : P3808-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
X0B24

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP021816.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.022	4	6	28	rVB	8922727	11045109	51.35%	3.579%
2	3.275	45	49	55	rVB	72020	84513	0.39%	0.027%
3	3.611	100	106	114	rBV	2037461	2711949	12.61%	0.879%
4	3.681	114	118	136	rVB	909702	1356534	6.31%	0.440%
5	4.011	170	174	182	rVB	60788	79570	0.37%	0.026%
6	4.093	185	188	205	rVB	57681	106662	0.50%	0.035%
7	4.828	304	313	320	rVB	89687	135645	0.63%	0.044%
8	4.916	323	328	344	rBV2	38211	104626	0.49%	0.034%
9	6.052	517	521	528	rBV6	16129	28149	0.13%	0.009%
10	6.952	666	674	684	rBV	1156818	1960278	9.11%	0.635%
11	7.110	694	701	708	rVB	919552	1466938	6.82%	0.475%
12	7.199	709	716	728	rVB	2266165	3410186	15.86%	1.105%
13	7.310	728	735	745	rBV	1083414	1816818	8.45%	0.589%
14	7.663	787	795	802	rBV	1461092	2254841	10.48%	0.731%
15	7.775	805	814	828	rVB	932326	1501611	6.98%	0.487%
16	8.122	866	873	883	rBV	1842155	3072694	14.29%	0.996%
17	8.475	926	933	942	rVV	1416493	2454006	11.41%	0.795%
18	8.575	943	950	961	rVB	862543	1419911	6.60%	0.460%
19	8.852	988	997	1003	rBV	2001934	3307932	15.38%	1.072%
20	8.922	1003	1009	1012	rVV	1028109	1805810	8.40%	0.585%
21	8.963	1012	1016	1026	rVB	2159745	3544878	16.48%	1.149%
22	9.493	1093	1106	1122	rBV	6678000	11066181	51.45%	3.586%
23	9.640	1122	1131	1145	rVV	844650	1446241	6.72%	0.469%
24	9.793	1152	1157	1166	rVB2	77620	133629	0.62%	0.043%
25	9.969	1176	1187	1201	rVB	5526271	8709529	40.49%	2.822%
26	10.187	1216	1224	1238	rBV	1413388	2335460	10.86%	0.757%
27	10.304	1240	1244	1253	rVB3	21131	38669	0.18%	0.013%
28	10.428	1255	1265	1273	rBV	4796924	8197336	38.11%	2.656%
29	10.557	1278	1287	1291	rBV	1209558	2235552	10.39%	0.724%
30	10.610	1291	1296	1303	rVV	6782457	11198055	52.06%	3.628%
31	10.687	1303	1309	1321	rVB	1240035	2348628	10.92%	0.761%
32	10.898	1337	1345	1360	rVB	6181338	9951168	46.27%	3.224%
33	11.234	1395	1402	1407	rBV	34707	65050	0.30%	0.021%
34	11.304	1407	1414	1420	rVB5	22644	41937	0.19%	0.014%
35	11.781	1491	1495	1504	rVB	33495	57632	0.27%	0.019%
36	12.045	1534	1540	1547	rBV6	24408	59074	0.27%	0.019%

LSC Area Percent Report

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Acq On : 04 Sep 2024 12:38
Operator : RC/JU
Sample : P3808-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
X0B24

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Title : SVOA CALIBRATION

37	12.110	1547	1551	1553	rVV	18170	27081	0.13%	0.009%
38	12.140	1553	1556	1561	rVV	24517	41025	0.19%	0.013%
39	12.216	1561	1569	1583	rVB	4232360	6910292	32.13%	2.239%
40	12.345	1583	1591	1601	rBV2	113080	230394	1.07%	0.075%
41	12.434	1601	1606	1614	rVB	62791	110678	0.51%	0.036%
42	12.569	1621	1629	1636	rBV	3066094	5023432	23.36%	1.628%
43	12.634	1636	1640	1646	rVB2	39955	65998	0.31%	0.021%
44	13.275	1737	1749	1761	rBV	6083104	9733315	45.25%	3.154%
45	13.451	1774	1779	1784	rBV	104216	148873	0.69%	0.048%
46	13.516	1784	1790	1795	rVV3	43037	66290	0.31%	0.021%
47	13.563	1795	1798	1807	rVB	34969	57166	0.27%	0.019%
48	13.810	1832	1840	1844	rBV	2361251	3440233	15.99%	1.115%
49	13.857	1844	1848	1855	rVB	3303306	4479869	20.83%	1.452%
50	13.975	1862	1868	1878	rVV	836884	1218248	5.66%	0.395%
51	14.098	1881	1889	1891	rBV2	2655707	4143970	19.27%	1.343%
52	14.128	1891	1894	1908	rVB	2988151	4293070	19.96%	1.391%
53	14.410	1935	1942	1951	rBV	1824917	2738508	12.73%	0.887%
54	14.616	1969	1977	1992	rBV	1149270	1913821	8.90%	0.620%
55	14.734	1992	1997	2002	rVV2	55116	112520	0.52%	0.036%
56	14.822	2006	2012	2025	rVB3	40048	104109	0.48%	0.034%
57	15.245	2077	2084	2092	rBV	3623535	5627608	26.16%	1.823%
58	15.422	2108	2114	2118	rBV	3804095	5125551	23.83%	1.661%
59	15.475	2118	2123	2129	rVV	14351274	21508462	100.00%	6.969%
60	15.539	2129	2134	2145	rVB	1009379	1459041	6.78%	0.473%
61	15.669	2154	2156	2163	rVB4	17300	23413	0.11%	0.008%
62	15.734	2163	2167	2173	rVB	60116	78159	0.36%	0.025%
63	15.963	2202	2206	2212	rBV4	27925	44418	0.21%	0.014%
64	16.086	2222	2227	2233	rVV2	68882	96666	0.45%	0.031%
65	16.398	2267	2280	2290	rBV	5630731	8354989	38.85%	2.707%
66	16.510	2292	2299	2312	rBV	5717486	8377215	38.95%	2.714%
67	17.010	2378	2384	2386	rBV6	27289	48367	0.22%	0.016%
68	17.033	2386	2388	2393	rVB	27783	34935	0.16%	0.011%
69	17.222	2413	2420	2423	rBV	2328890	3447871	16.03%	1.117%
70	17.263	2423	2427	2432	rVV	4258746	6368011	29.61%	2.063%
71	17.328	2432	2438	2440	rVV	3649276	5565983	25.88%	1.803%
72	17.363	2440	2444	2461	rVV	4808263	7531693	35.02%	2.440%
73	17.504	2464	2468	2475	rVB2	19824	32382	0.15%	0.010%
74	17.639	2487	2491	2497	rVB	32444	46704	0.22%	0.015%
75	17.857	2524	2528	2531	rBV2	22001	32772	0.15%	0.011%
76	18.163	2576	2580	2586	rBV3	39448	70298	0.33%	0.023%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
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Acq On : 04 Sep 2024 12:38
Operator : RC/JU
Sample : P3808-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
X0B24

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Title : SVOA CALIBRATION

77	18.233	2586	2592	2604	rVV	5877211	7894723	36.71%	2.558%
78	18.686	2664	2669	2677	rVB	257357	402556	1.87%	0.130%
79	19.251	2760	2765	2769	rBV2	45249	80577	0.37%	0.026%
80	19.322	2773	2777	2780	rVV	48411	64427	0.30%	0.021%
81	19.698	2835	2841	2850	rBV	5550469	7150734	33.25%	2.317%
82	21.674	3172	3177	3181	rBV2	2619547	4160618	19.34%	1.348%
83	21.727	3181	3186	3194	rVB	5436605	8844312	41.12%	2.866%
84	22.763	3358	3362	3370	rVB	325849	575823	2.68%	0.187%
85	22.904	3379	3386	3394	rBV	2833429	5086577	23.65%	1.648%
86	24.004	3564	3573	3579	rBV	2947970	7273469	33.82%	2.357%
87	24.080	3579	3586	3597	rVB	6012354	13389681	62.25%	4.338%
88	24.845	3707	3716	3722	rBV	2970142	7713235	35.86%	2.499%
89	24.915	3722	3728	3744	rVB	3285027	8409476	39.10%	2.725%
90	25.080	3747	3756	3769	rVB	1607544	4528413	21.05%	1.467%
91	28.956	4400	4415	4423	rBV	2777326	12765191	59.35%	4.136%

Sum of corrected areas: 308626043

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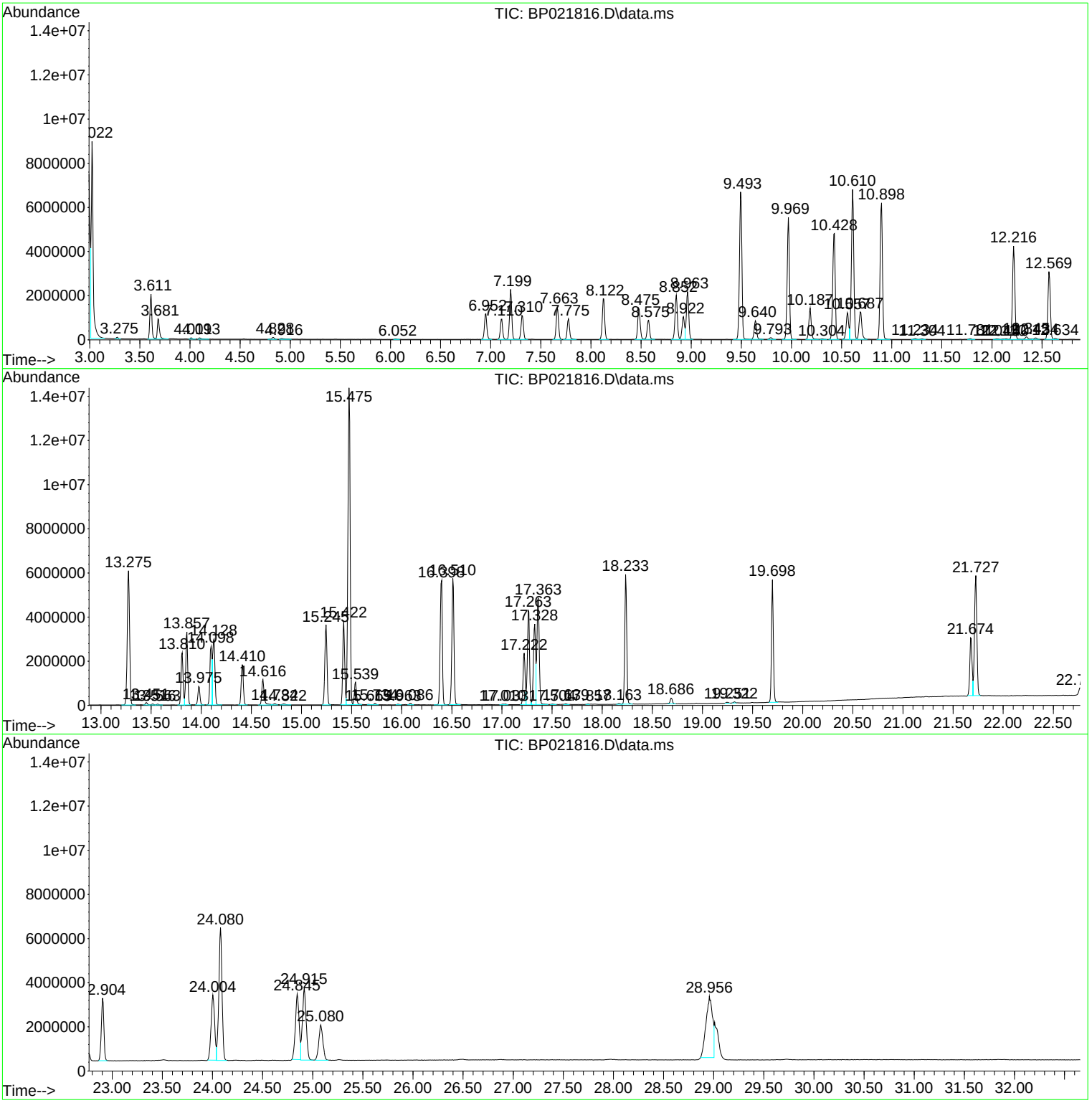
Operator : RC/JU
Sample : P3808-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
X0B24

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
Data File : BP021816.D
Acq On : 04 Sep 2024 12:38
Operator : RC/JU
Sample : P3808-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
X0B24

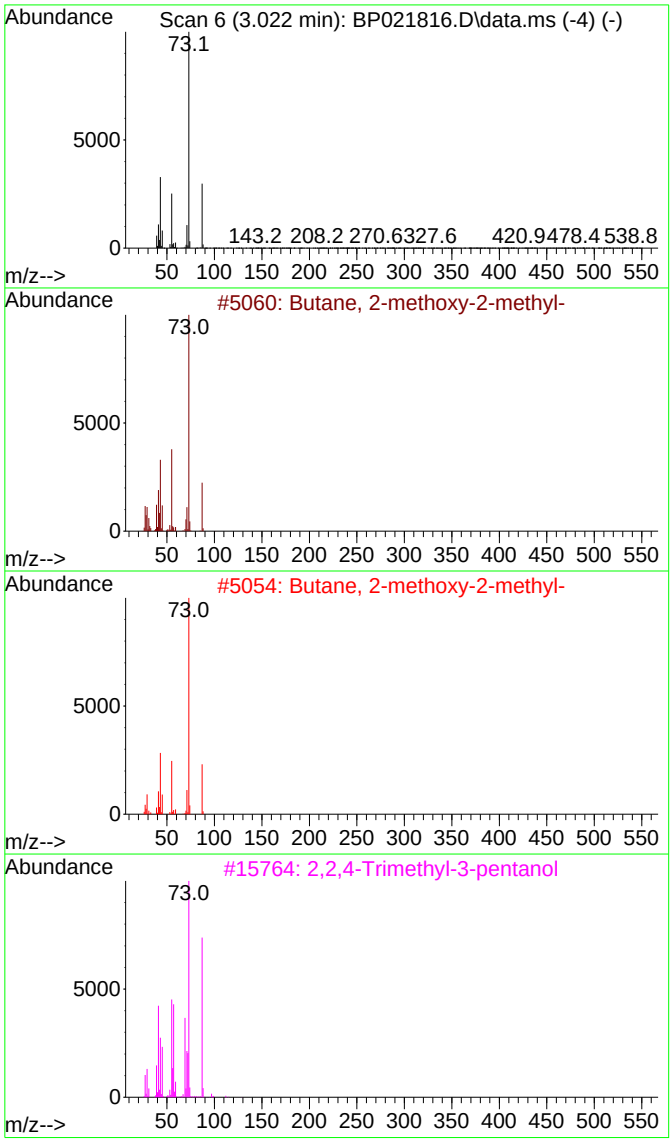
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

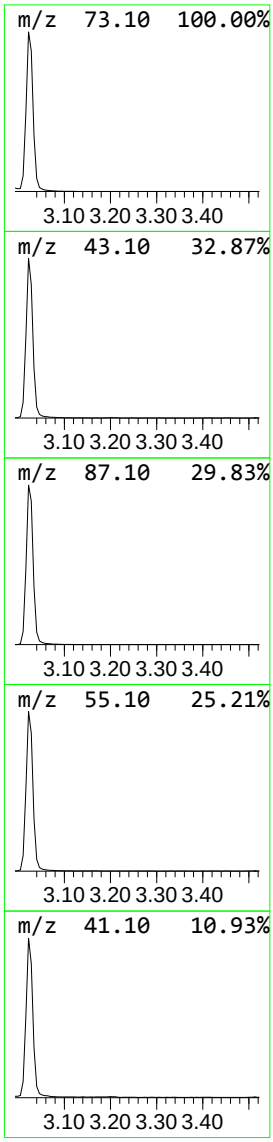
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.022	147.11 ng/ul	11045100	1,4-Dichlorobenzene-d4	7.775

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	59
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
5	N-Ethylformamide	73	C3H7NO	000627-45-2	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
Data File : BP021816.D
Acq On : 04 Sep 2024 12:38
Operator : RC/JU
Sample : P3808-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
X0B24

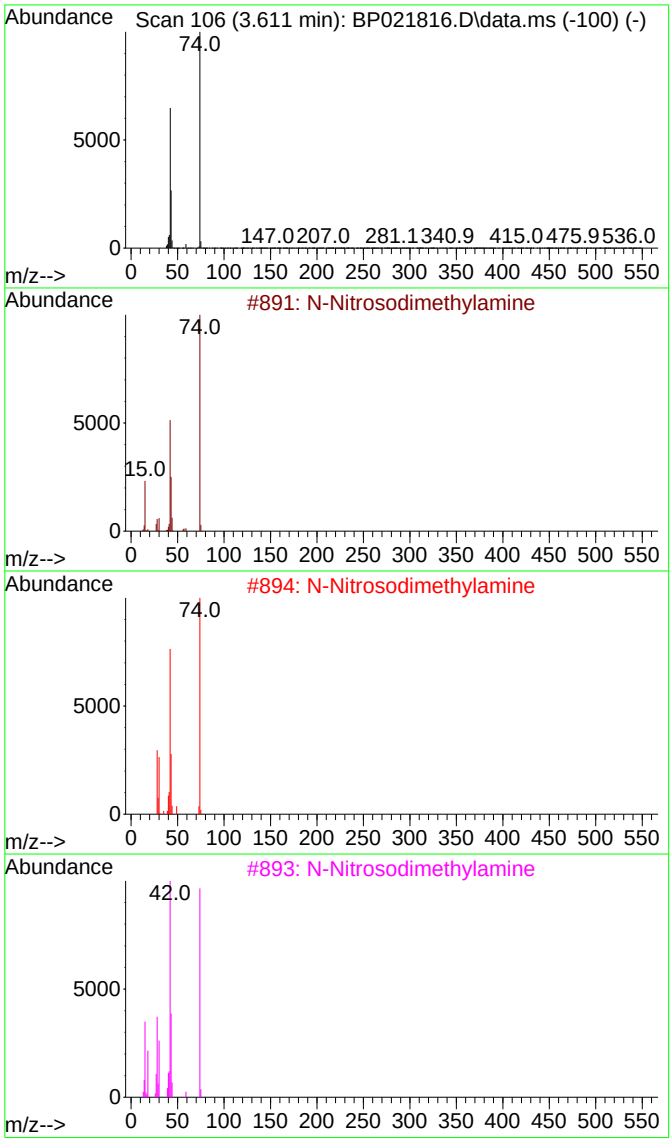
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

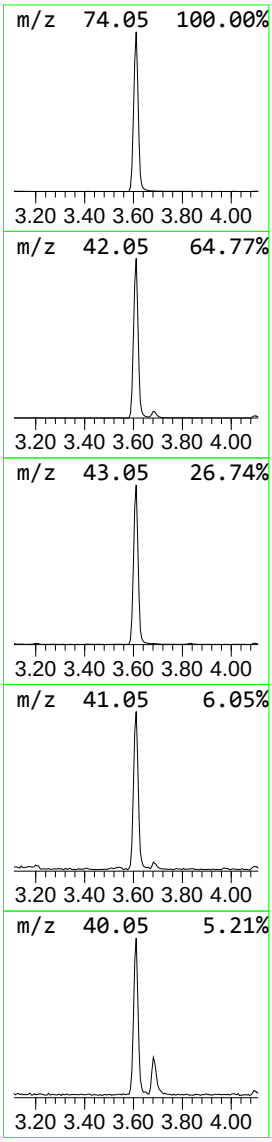
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 N-Nitrosodimethylamine Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.611	36.12 ng/ul	2711950	1,4-Dichlorobenzene-d4	7.775

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	N-Nitrosodimethylamine	74	C2H6N2O	000062-75-9	78
2	N-Nitrosodimethylamine	74	C2H6N2O	000062-75-9	9
3	N-Nitrosodimethylamine	74	C2H6N2O	000062-75-9	5
4	N-Nitrosodimethylamine	74	C2H6N2O	000062-75-9	4
5	Aminoguanidine	74	CH6N4	000079-17-4	4





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
Data File : BP021816.D
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Operator : RC/JU
Sample : P3808-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
X0B24

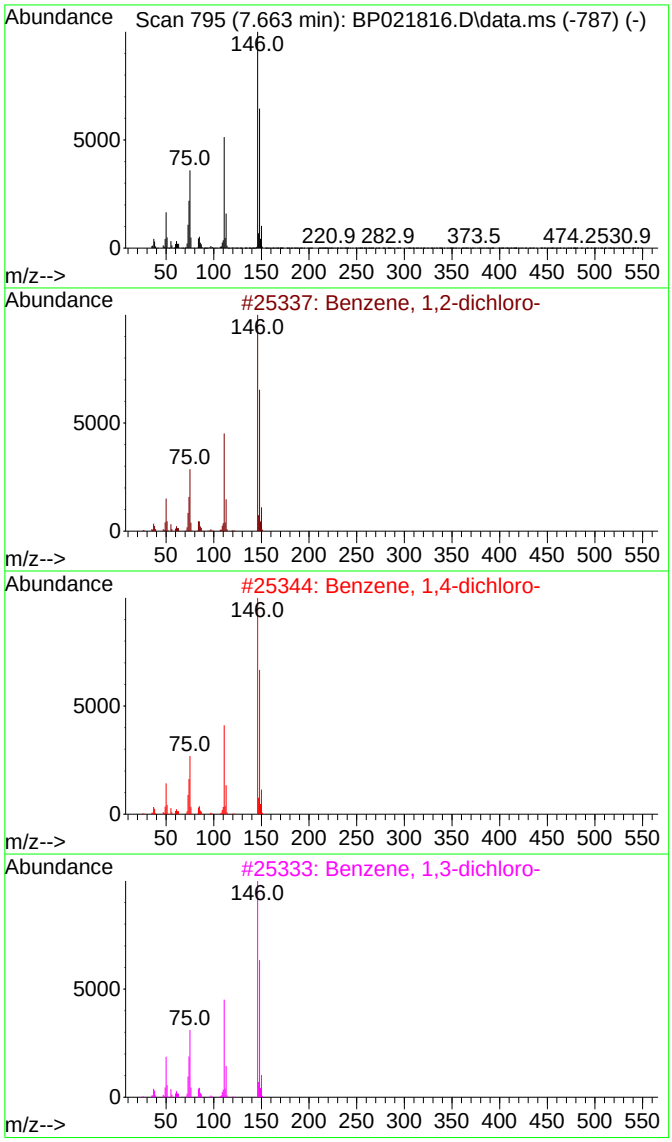
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

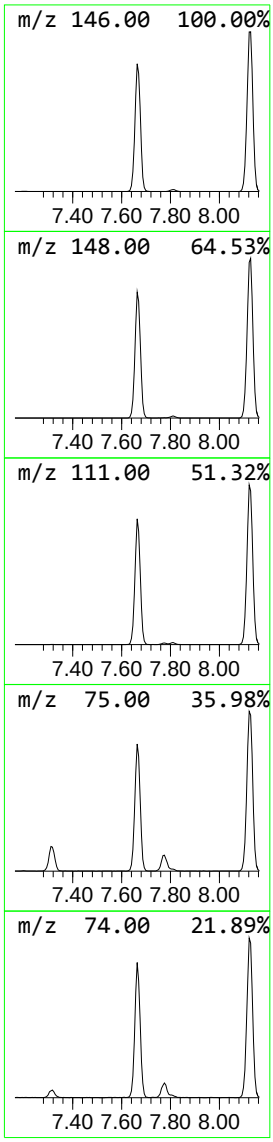
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1,2-dichloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.663	30.03 ng/ul	2254840	1,4-Dichlorobenzene-d4	7.775

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
2	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
3	Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
4	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	97
5	Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
Data File : BP021816.D
Acq On : 04 Sep 2024 12:38
Operator : RC/JU
Sample : P3808-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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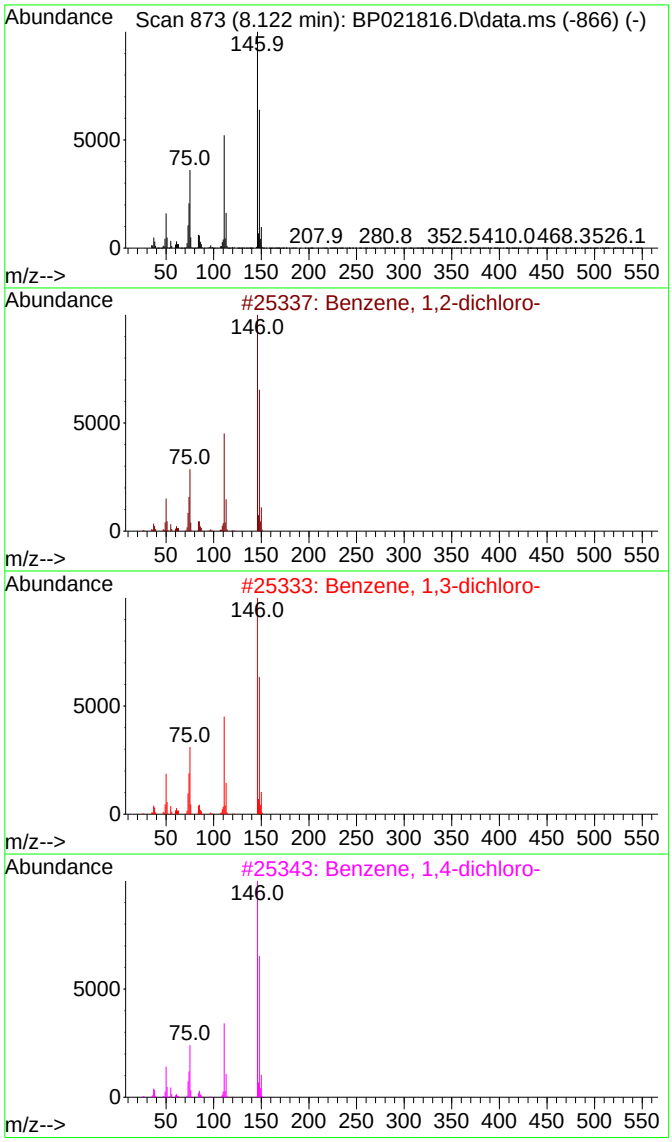
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

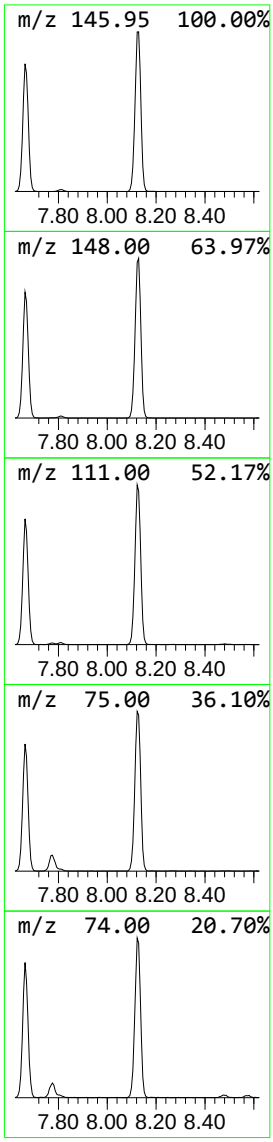
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1,3-dichloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.122	40.93 ng/ul	3072690	1,4-Dichlorobenzene-d4	7.775

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
2	Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
3	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	97
4	Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	96
5	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
Data File : BP021816.D
Acq On : 04 Sep 2024 12:38
Operator : RC/JU
Sample : P3808-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
X0B24

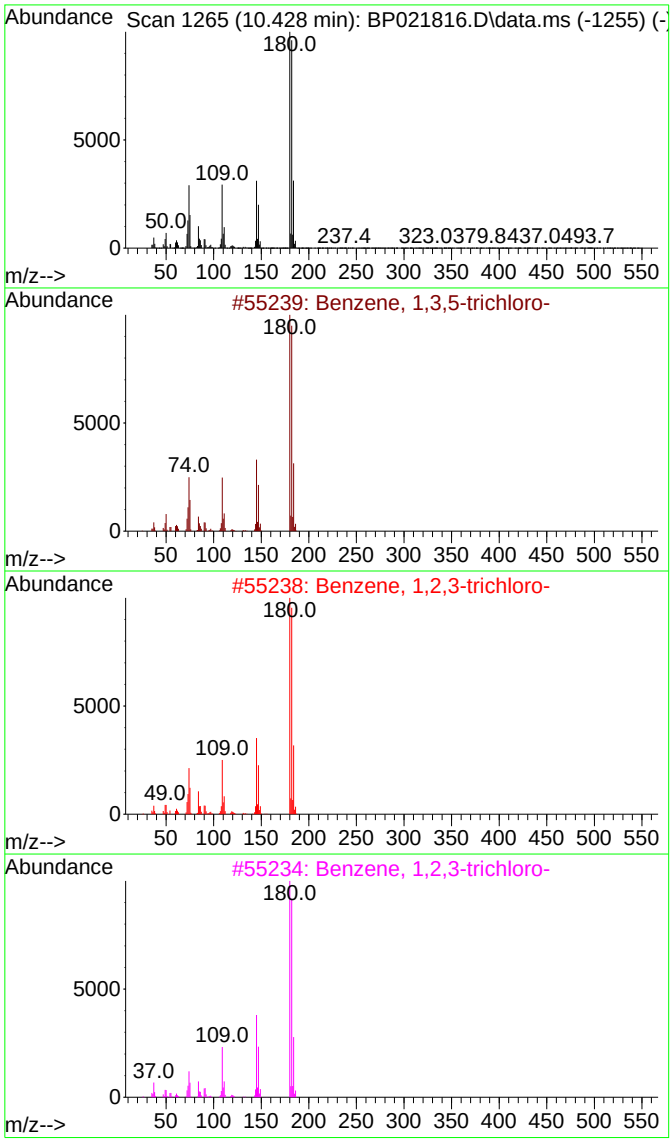
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

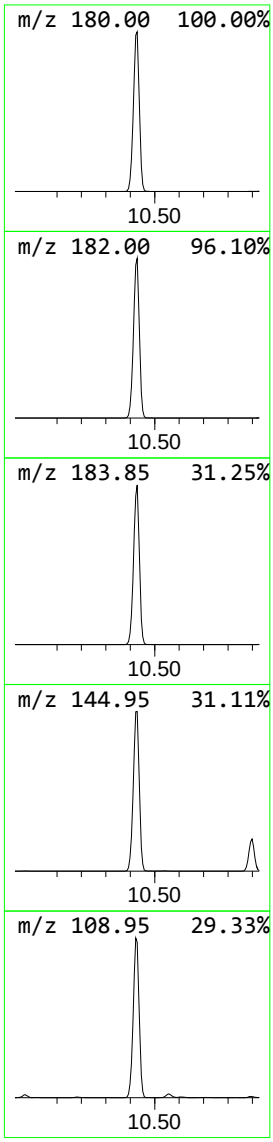
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1,3,5-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.428	73.34 ng/ul	8197340	Naphthalene-d8	10.557

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
2	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
3	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
4	Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	97
5	Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	97





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
Data File : BP021816.D
Acq On : 04 Sep 2024 12:38
Operator : RC/JU
Sample : P3808-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
X0B24

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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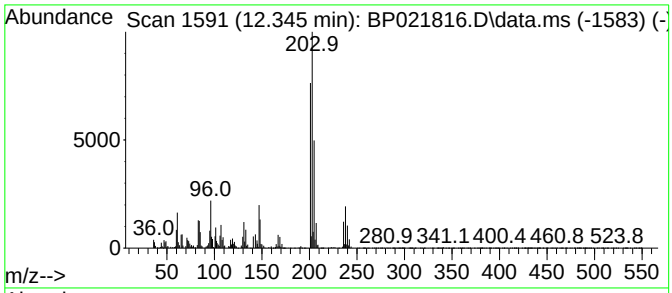
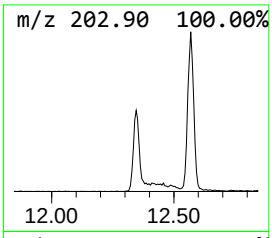
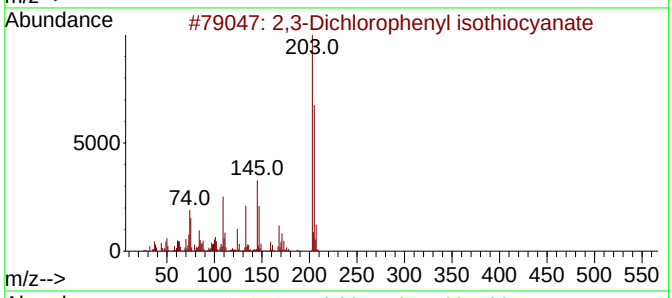
Peak Number 6 unknown-01 Concentration Rank 8

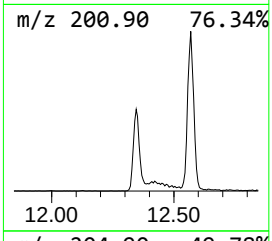
R.T.	EstConc	Area	Relative to ISTD	R.T.
12.345	2.06 ng/ul	230394	Naphthalene-d8	10.557

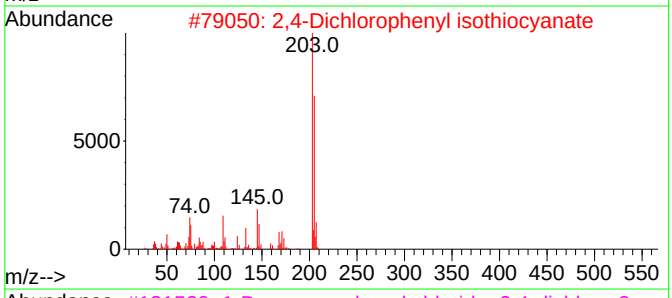
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3-Dichlorophenyl isothiocyanate	203	C7H3Cl2NS	006590-97-2	45
2	2,4-Dichlorophenyl isothiocyanate	203	C7H3Cl2NS	006590-96-1	38
3	1-Benzenecarbonyl chloride, 2,4-...	238	C8H5Cl3O2	1000196-85-1	32
4	2,6-Dichloro-4-(1,1-dimethylethy...	218	C10H12Cl2O	1000305-55-5	25
5	3,4-Dichlorophenyl isothiocyanate	203	C7H3Cl2NS	006590-94-9	25

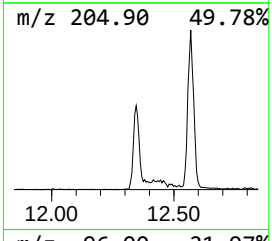


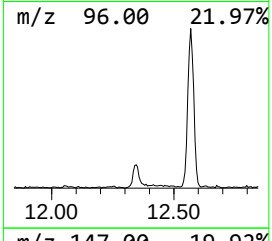


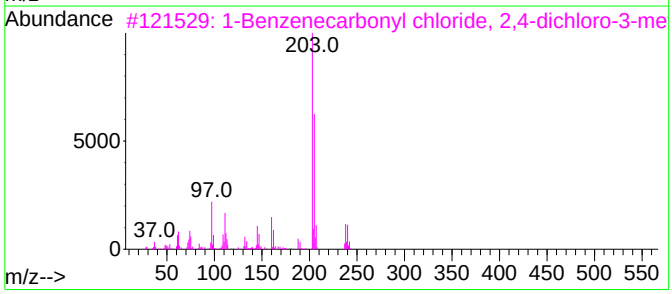


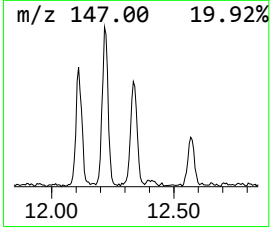












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
Data File : BP021816.D
Acq On : 04 Sep 2024 12:38
Operator : RC/JU
Sample : P3808-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
X0B24

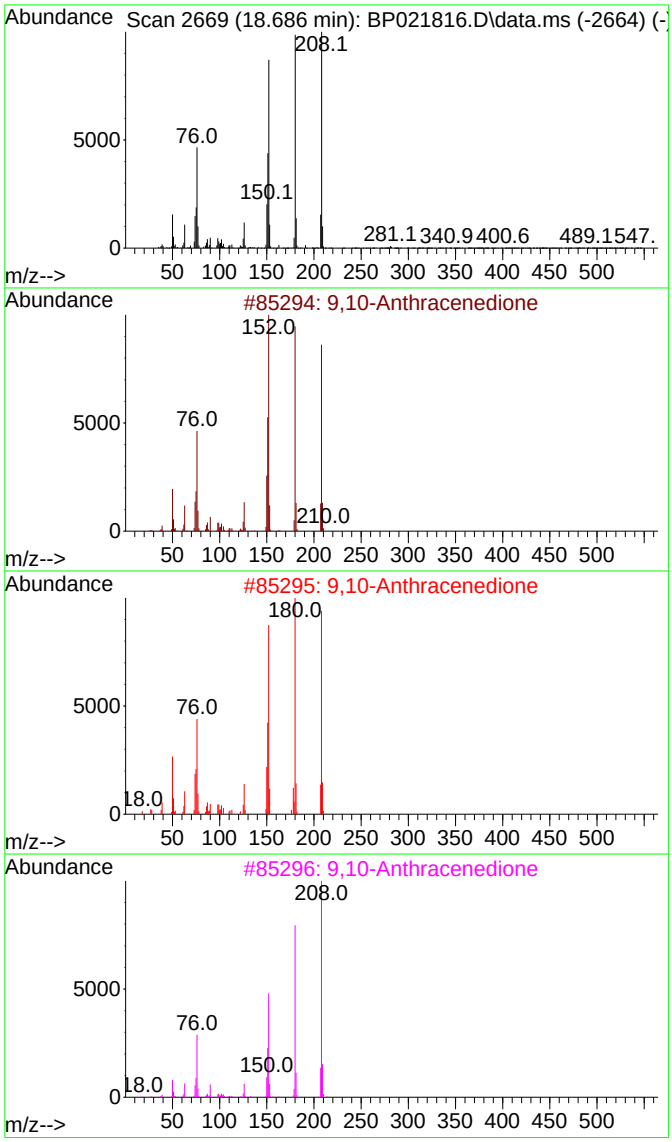
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

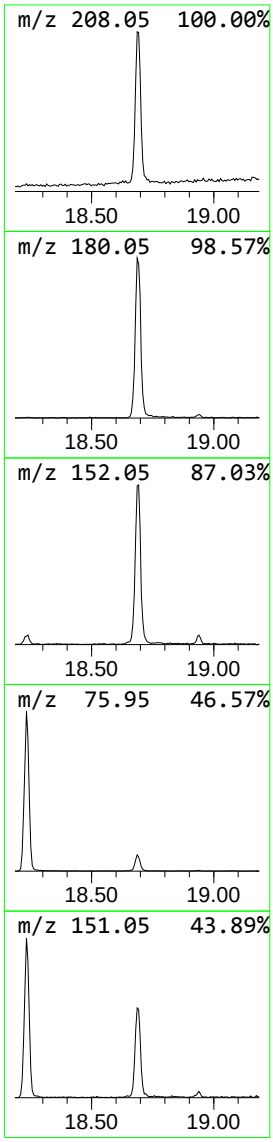
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 9,10-Anthracenedione Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.686	2.34 ng/ul	402556	Phenanthrene-d10	17.222

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
2	9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
3	9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
4	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	91
5	9,10-Anthracenedione	208	C14H8O2	000084-65-1	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
Data File : BP021816.D
Acq On : 04 Sep 2024 12:38
Operator : RC/JU
Sample : P3808-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
X0B24

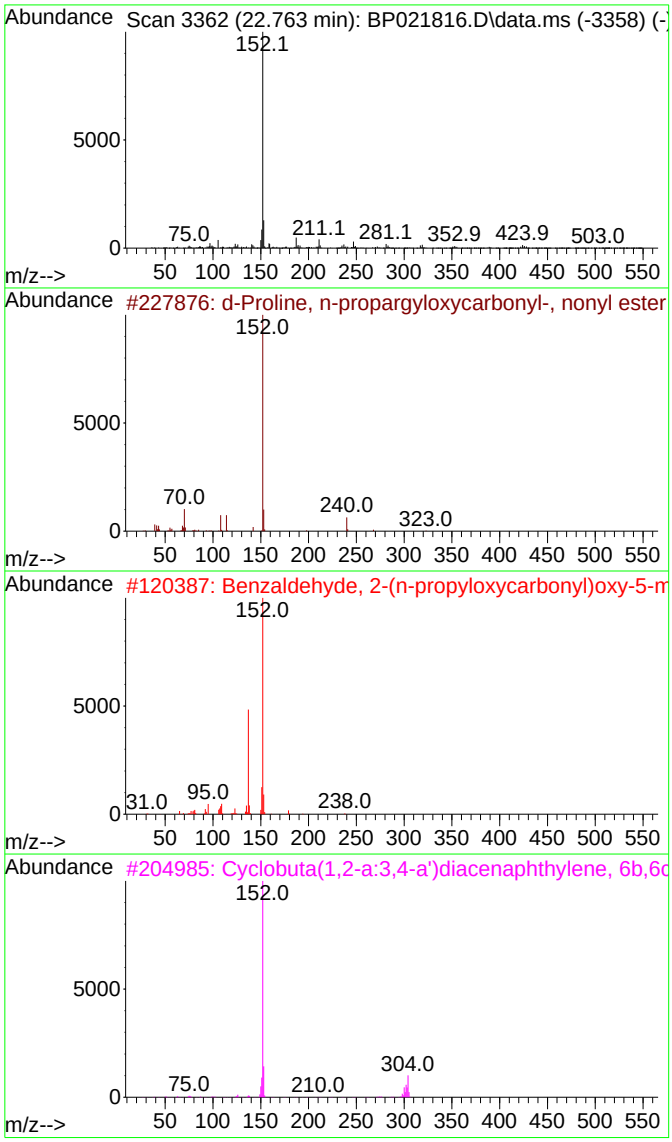
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

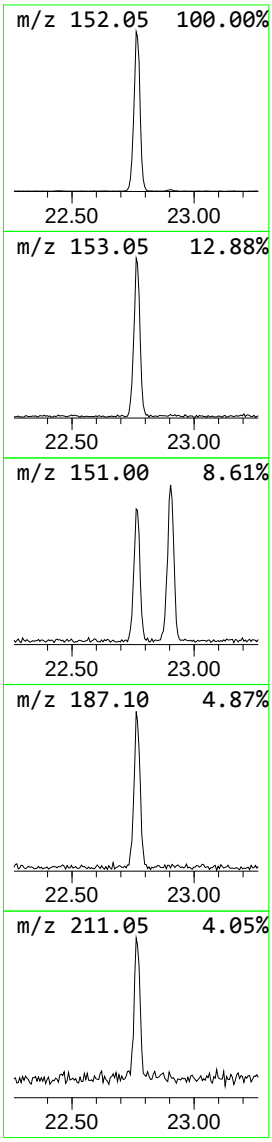
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 d-Proline, n-propargyloxyca... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.763	2.77 ng/ul	575823	Chrysene-d12	21.674

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	d-Proline, n-propargyloxycarbony...	323	C18H29NO4	1000328-07-1	64
2	Benzaldehyde, 2-(n-propyloxycarb...	238	C12H14O5	1000487-65-6	64
3	Cyclobuta(1,2-a:3,4-a')diacenaph...	304	C24H16	007259-03-2	64
4	Acenaphthylene	152	C12H8	000208-96-8	64
5	Phthalic acid, ethyl 2H-octahydr...	345	C20H27NO4	1000322-80-1	50





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
Data File : BP021816.D
Acq On : 04 Sep 2024 12:38
Operator : RC/JU
Sample : P3808-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
X0B24

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-metho...	3.022	147.1	ng/ul	11045100	1	7.775	1501610	20.0
N-Nitrosodimeth...	3.611	36.1	ng/ul	2711950	1	7.775	1501610	20.0
Benzene, 1,2-di...	7.663	30.0	ng/ul	2254840	1	7.775	1501610	20.0
Benzene, 1,3-di...	8.122	40.9	ng/ul	3072690	1	7.775	1501610	20.0
Benzene, 1,3,5-...	10.428	73.3	ng/ul	8197340	2	10.557	2235550	20.0
unknown-01	12.345	2.1	ng/ul	230394	2	10.557	2235550	20.0
9,10-Anthracene...	18.686	2.3	ng/ul	402556	4	17.222	3447870	20.0
d-Proline, n-pr...	22.763	2.8	ng/ul	575823	5	21.674	4160620	20.0